

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078530.D
 Acq On : 15 Apr 2015 8:13
 Operator : TP/IZ
 Sample : G1823-05DL 10X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3DL

Quant Time: Apr 15 11:49:20 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 01:10:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	24310	20.00	ng	0.00
21) Naphthalene-d8	8.39	136	104184	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	55442	20.00	ng	0.00
63) Phenanthrene-d10	12.37	188	116704	20.00	ng	0.00
75) Chrysene-d12	15.62	240	137776	20.00	ng	-0.03
86) Perylene-d12	17.37	264	139597	20.00	ng	-0.11
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	13612	9.23	ng	0.01
7) Phenol-d6	6.42	99	18694	10.12	ng	0.01
23) Nitrobenzene-d5	7.51	82	10644	6.19	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	5040	10.96	ng	0.00
44) 2-Fluorobiphenyl	9.74	172	27318	7.04	ng	0.00
78) Terphenyl-d14	14.35	244	37468	6.15	ng	0.00
Target Compounds						Qvalue
51) Acenaphthene	10.58	154	12813	3.95	ng	97
57) Fluorene	11.21	166	10119	2.52	ng	99
70) Phenanthrene	12.39	178	170195	25.21	ng	99
71) Anthracene	12.46	178	47833	7.19	ng	100
72) Carbazole	12.67	167	18092	2.90	ng	97
74) Fluoranthene	13.86	202	349586	49.10	ng	92
77) Pyrene	14.14	202	330208	38.18	ng	98
80) Benzo(a)anthracene	15.61	228	194947	23.78	ng	98
82) Chrysene	15.66	228	166278	21.59	ng	99
85) Indeno(1,2,3-cd)pyrene	18.62	276	94616m	10.83	ng	
87) Benzo(b)fluoranthene	16.93	252	207672m	24.07	ng	
88) Benzo(k)fluoranthene	16.94	252	91988m	12.00	ng	
89) Benzo(a)pyrene	17.30	252	149703	19.88	ng	# 94
90) Dibenzo(a,h)anthracene	18.64	278	24625m	3.27	ng	
91) Benzo(g,h,i)perylene	18.97	276	88493m	11.71	ng	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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